

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112624\
 Data File : BN035348.D
 Acq On : 27 Nov 2024 13:34
 Operator : RC/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4356

Quant Time: Nov 27 23:49:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 06:09:15 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.305	152	3418	0.400	ng/ul	0.00
4) Naphthalene-d8	10.053	136	8838	0.400	ng/ul	-0.01
9) Acenaphthene-d10	13.967	164	5286	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.734	188	10462	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.976	240	6187	0.400	ng/ul	# 0.00
23) Perylene-d12	23.071	264	6399	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.972	96	1352	0.433	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.659	152	4822	0.372	ng/ul	0.00
18) Fluoranthene-d10	18.786	212	8450	0.495	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.001	88	1431	0.422	ng/ul	96
5) Naphthalene	10.102	128	9180	0.412	ng/ul	99
7) 2-Methylnaphthalene	11.730	142	6170	0.400	ng/ul	99
8) 1-Methylnaphthalene	11.956	142	6253	0.401	ng/ul	99
10) Acenaphthylene	13.676	152	9058	0.432	ng/ul	99
11) Acenaphthene	14.027	153	6943	0.433	ng/ul	98
12) Fluorene	15.031	166	7701	0.409	ng/ul	99
14) Pentachlorophenol	16.392	266	790	0.324	ng/ul	98
15) Phenanthrene	16.776	178	11745	0.425	ng/ul	99
16) Anthracene	16.865	178	10211	0.398	ng/ul	100
19) Fluoranthene	18.814	202	11407	0.530	ng/ul#	93
20) Pyrene	19.181	202	11300	0.504	ng/ul#	91
21) Benzo(a)anthracene	20.961	228	8474	0.395	ng/ul	99
22) Chrysene	21.014	228	9385	0.434	ng/ul	100
24) Benzo(b)fluoranthene	22.457	252	8877	0.420	ng/ul	95
25) Benzo(k)fluoranthene	22.498	252	9706	0.445	ng/ul#	96
26) Benzo(a)pyrene	22.981	252	8079	0.418	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	25.111	276	10823	0.415	ng/ul#	77
28) Dibenzo(a,h)anthracene	25.128	278	8178	0.400	ng/ul#	95
29) Benzo(g,h,i)perylene	25.725	276	8903	0.433	ng/ul#	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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